

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012320\
 Data File : BM024611.D
 Acq On : 23 Jan 2020 17:09
 Operator : CG/JU
 Sample : K5111-07
 Misc : 5 PPM LOD
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 LOD-MDL-WATER-01-QT4-2019

Manual Integrations
 APPROVED

mohammad
 1/24/2020 1:55:13 PM

Quant Time: Jan 24 11:45:35 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM012320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 23 14:57:07 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.45	152	199214	20.00	ng	0.00
21) Naphthalene-d8	10.20	136	815696	20.00	ng	0.00
39) Acenaphthene-d10	14.09	164	585422	20.00	ng	0.00
64) Phenanthrene-d10	16.84	188	1384558	20.00	ng	0.00
76) Chrysene-d12	21.05	240	1488438	20.00	ng	0.00
87) Perylene-d12	23.17	264	1675975	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.08	112	1354392	129.29	ng	0.00
7) Phenol-d6	6.64	99	1769848	122.65	ng	0.00
23) Nitrobenzene-d5	8.58	82	1310649	78.80	ng	0.00
42) 2,4,6-Tribromophenol	15.59	330	1255697	131.86	ng	0.00
45) 2-Fluorobiphenyl	12.71	172	3424587	79.53	ng	0.00
79) Terphenyl-d14	19.50	244	7021025	88.06	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.08	88	21642	5.223	ng	96
3) Pyridine	3.46	79	58695	4.886	ng	# 92
4) n-Nitrosodimethylamine	3.37	42	25086	4.705	ng	# 98
6) Aniline	6.79	93	82056	4.671	ng	98
8) 2-Chlorophenol	7.02	128	55619	4.664	ng	94
9) Benzaldehyde	6.60	77	40146m	4.708	ng	
10) Phenol	6.66	94	66926	4.653	ng	83
11) bis(2-Chloroethyl)ether	6.89	93	56062	4.868	ng	95
12) 1,3-Dichlorobenzene	7.33	146	70386	4.805	ng	97
13) 1,4-Dichlorobenzene	7.48	146	72615	4.886	ng	94
14) 1,2-Dichlorobenzene	7.79	146	69630	4.854	ng	96
15) Benzyl Alcohol	7.69	79	53692	4.513	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.97	45	49849	4.729	ng	96
17) 2-Methylphenol	7.89	107	44260	4.366	ng	94
18) Hexachloroethane	8.51	117	28324	4.935	ng	98
19) n-Nitroso-di-n-propylamine	8.25	70	49696	4.587	ng	97
20) 3+4-Methylphenols	8.22	107	61491	4.382	ng	98
22) Acetophenone	8.26	105	100353	4.771	ng	# 98
24) Nitrobenzene	8.62	77	80821m	5.071	ng	
25) Isophorone	9.15	82	130525	4.618	ng	99
26) 2-Nitrophenol	9.33	139	28656	3.883	ng	97
27) 2,4-Dimethylphenol	9.41	122	43908	4.378	ng	94
28) bis(2-Chloroethoxy)methane	9.64	93	79267	4.591	ng	98
29) 2,4-Dichlorophenol	9.86	162	55382	4.087	ng	99
30) 1,2,4-Trichlorobenzene	10.07	180	76151	4.660	ng	99
31) Naphthalene	10.25	128	192259	4.667	ng	99
32) Benzoic acid	9.48	122	29427m	3.207	ng	
33) 4-Chloroaniline	10.37	127	80950	4.483	ng	99
34) Hexachlorobutadiene	10.56	225	56516	5.003	ng	93
35) Caprolactam	11.13	113	16617	3.822	ng	91
36) 4-Chloro-3-methylphenol	11.50	107	57266	3.975	ng	98
37) 2-Methylnaphthalene	11.87	142	147462	4.638	ng	98
38) 1-Methylnaphthalene	12.10	142	142159	4.701	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.26	216	96309	4.609	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.24	237	47452	4.052	ng	96
43) 2,4,6-Trichlorophenol	12.50	196	53897	4.054	ng	97
44) 2,4,5-Trichlorophenol	12.57	196	55012	4.051	ng	95
46) 1,1'-Biphenyl	12.92	154	212435	4.795	ng	99
47) 2-Chloronaphthalene	12.95	162	165712	4.633	ng	99
48) 2-Nitroaniline	13.16	65	42789	3.923	ng	96
49) Acenaphthylene	13.80	152	242055	4.521	ng	99
50) Dimethylphthalate	13.56	163	219523	4.637	ng	98
51) 2,6-Dinitrotoluene	13.67	165	42180	4.191	ng	88
52) Acenaphthene	14.16	154	150264	4.523	ng	99
53) 3-Nitroaniline	14.00	138	40067	3.838	ng	94
54) 2,4-Dinitrophenol	14.21	184	16210	2.693	ng	94
55) Dibenzofuran	14.49	168	254285	4.693	ng	98
56) 4-Nitrophenol	14.32	139	28660	3.512	ng	98
57) 2,4-Dinitrotoluene	14.46	165	57302	3.973	ng	93
58) Fluorene	15.14	166	202260	4.442	ng	97
59) 2,3,4,6-Tetrachlorophenol	14.73	232	57761	4.195	ng	96
60) Diethylphthalate	14.94	149	215458	4.483	ng	99
61) 4-Chlorophenyl-phenylether	15.16	204	120291	4.582	ng	97
62) 4-Nitroaniline	15.16	138	42015	3.646	ng	86
63) Azobenzene	15.44	77	183829	4.484	ng	98
65) 4,6-Dinitro-2-methylphenol	15.23	198	25203	3.096	ng	96
66) n-Nitrosodiphenylamine	15.37	169	176438	4.436	ng	98
67) 4-Bromophenyl-phenylether	16.05	248	75769	4.401	ng	97
68) Hexachlorobenzene	16.16	284	93652	4.740	ng	96
69) Atrazine	16.33	200	64828	4.246	ng	98
70) Pentachlorophenol	16.50	266	43666	3.628	ng	98
71) Phenanthrene	16.88	178	343073	4.576	ng	100
72) Anthracene	16.97	178	319573	4.356	ng	99
73) Carbazole	17.24	167	286536	4.290	ng	98
74) Di-n-butylphthalate	17.84	149	343992	4.029	ng	100
75) Fluoranthene	18.91	202	410540	4.387	ng	98
77) Benzidine	19.11	184	119789	3.456	ng	99
78) Pyrene	19.27	202	431429	4.397	ng	100
80) Butylbenzylphthalate	20.22	149	150457	3.654	ng	100
81) Benzo(a)anthracene	21.04	228	470585	4.555	ng	98
82) 3,3'-Dichlorobenzidine	20.98	252	144697	3.894	ng	100
83) Chrysene	21.09	228	455413	4.620	ng	100
84) Bis(2-ethylhexyl)phthalate	21.01	149	226615	3.730	ng	98
85) Di-n-octyl phthalate	21.86	149	363909	3.626	ng	99
86) Indeno(1,2,3-cd)pyrene	25.23	276	478865	4.457	ng	99
88) Benzo(b)fluoranthene	22.54	252	452239	4.160	ng	99
89) Benzo(k)fluoranthene	22.59	252	466926	4.401	ng	99
90) Benzo(a)pyrene	23.07	252	413320	4.182	ng	99
91) Dibenzo(a,h)anthracene	25.24	278	404019	4.307	ng	98
92) Benzo(g,h,i)perylene	25.85	276	367357	4.313	ng	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

